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BADEM			Application Number		10/825,186
			Filing Date		April 16, 2004
FORM (to be used for all correspondence after initial filing)		First Named Inventor		Zhang et al.	
			Group Art Unit		1645
			Confirmation No.		8260
Total Number of Pages in This Su	ubmission		Attorney Docket Number		057953-1221
		ENCLOSU	RES (check all that apply)		
Fee Transmittal Form Fee Attached Preliminary Amendment After Final Affidavits/declaration(s) Extension of Time Request for Express Abandonment Request Market Mark	s of ities Letter	Gor an A Submiss Drawing Declarat Licensir Petition Petition Applicat Power o Change Termina Request	tion and Power of Attorney ng-related Papers to Convert to a Provisional tion of Attorney, Revocation of Correspondence Address al Disclaimer for Refund mber of CD(s)	ayme	After Allowance Communication to Group Appeal Communication to Board of Appeals and Interferences Appeal Communication to Group (Appeal Notice, Brief, Reply Brief) Proprietary Information Status Letter Application Data Sheet Request for Corrected Filing Receipt with Enclosures A self-addressed prepaid postcard for acknowledging receipt Other Enclosure(s) (please identify below):
Firm or Individual name	aufman, Reg	14603-1051 1304	OR A	GENT	
Signature	•	1			
Date	December	1, 2004			
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(703) <u>872-9306</u>

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PATENT

Docket No.: 57953/1221 (ZHA01-01)

BADEN THE UNITED STATES PATENT AND TRADEMARK OFFICE

Applicant(s)	:	Zhang et al.) Examiner:) To Be Assigned
Serial No.	:	10/825,186)
Cnfrm. No.	:	8260) Art Unit:) 1645
Filed	:	April 16, 2004)
For	:	A METHOD FOR INTRODUCING CONJUGATED CAPS ONTO MOLECULAR FRAGMENTS AND SYSTEMS AND METHODS FOR USING THE SAME TO DETERMINE INTER-MOLECULAR INTERACTION ENERGIES)))))

INFORMATION DISCLOSURE STATEMENT UNDER 37 CFR §§ 1.97-1.98

Mail Stop: Amendment Commissioner for Patents P.O. Box 1450 Alexandria, VA 22313-1450

Dear Sir:

Pursuant to 37 CFR §§ 1.97-1.98, applicants hereby bring to the attention of the United States Patent and Trademark Office, the enclosed references listed on the attached PTO-1449 form.

Pursuant to 37 C.F.R. § 1.97(b)(3), no fee is required. If additional fees are required, however, the Commissioner is hereby authorized to charge any fees to Deposit Account No. 14-1138.

Respectfully submitted,

Date: 1/2004

Marc S. Kaufman Registration No. 35,212

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R813070.1

PTO/SB/08A (10-01)

Approved for use through 10/31/2002. OMB 0651-0031

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Substitute	for form 1449A/PTOR	ADEM		Complete if Known		
INFO	PMATION I	NSCI	AGUER	Application Number	10/825,186	
INFORMATION DISCLOSURE STATEMENT BY APPLICANT (use as many sheets as necessary)				Filing Date	April 16, 2004	
				First Named Inventor	Zhang et al.	
			(יכיי)	Art Unit	1645	
				Examiner Name	To Be Assigned	
Sheet	1	of	2	Attorney Docket Number	57953/1221 (ZHA01-01)	

			U.S. PATENT DOCUMEN	TS		
Examiner Cite Initials No.	U.S. Patent Document	Publication Date	Name of Patentee or	Pages, Columns, Lines, Where		
		Number - Kind Code ² (if known)	MM-DD-YYYY	Applicant of Cited Document	Relevant Passages or Relevant Figures Appear	
	1	US-				
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			FO	REIGN PATENT D	OCUMENTS				
Examiner Initials	Cite No. ¹	Foreign Patent Document Kind Code		Publication Date MM-DD-YYYY	Name of Patentee or Applicant of Cited Document	Pages, Columns, Lines, Where Relevant Passages or Relevant Figures Appear	т•		
		Country Code ³ Number ⁴	(if known)		177				
					 				
					1				
		ОТН	IER PRIOR A	ART – NON PATENT L	ITERATURE DOCUMENTS				
Examiner Initials*	Cite No.1	Include name of the author (in CAPITAL LETTERS), title of the article (when appropriate), title of the item (book, magazine, journal, serial, symposium, catalog, etc.)., date, page(s), volume-issue number(s), publisher, city and/or country where published.							
	1	Chen et al., "Fractionation of Peptide with Disulfide Bond for Quantum Mechanical Calculation of Interaction Energy with Molecules," <i>Journal of Chemical Physics</i> 120(2):839-844 (2004)							
	2	Chen et al., "Theoretical Method for Full ab initio Calculation of DNA/RNA-Ligand Interaction Energy," Journal of Chemical Physics 120(24):11386-11391 (2004)							
	Gao et al., "An Efficient Linear Scaling Method for ab initio Calculation of Electron Density of Proteins," Chemical Physics Letters 394:293-297 (2004)								

Examiner	Date	П
Signature	Considered	

^{*}EXAMINER: Initial if reference considered, whether or not citation is in conformance with MPEP 609. Draw line through citation if not in conformance and not considered. Include copy of this form with next communication to applicant.

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¹ Applicant's unique citation designation number (optional). ² See Kinds Codes of USPTO Patent Documents at www.uspto.gov or MPEP 901.04. ³ Enter Office that issued the document, by the two-letter code (WIPO Standard ST.3). ⁴ For Japanese patent documents, the indication of the year of the reign of the Emperor must precede the serial number of the patent document. ⁵ Kind of document by the appropriate symbols as indicated on the document under WIPO Standard ST.16 if possible. ⁶ Applicant is to place a check mark here if English language Translation is attached.

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PTO/SB/08B (10-01)

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Substitute for form 1449B/PTO RADENAM				Complete if Known		
INFO	INFORMATION DISCLOSURE			Application Number	10/825,186	
STATEMENT BY APPLICANT				Filing Date	April 16, 2004	
SIAI	(use as many sheets as necessary)			First Named Inventor	Zhang et al.	
				Group Art Unit	1645	
				Examiner Name	To Be Assigned	
Sheet	2	of	2	Attorney Docket Number	57953/1221 (ZHA01-01)	

		OTHER PRIOR ART – NON PATENT LITERATURE DOCUMENTS	
Examiner Initials*	Cite No.1	Include name of the author (in CAPITAL LETTERS), title of the article (when appropriate), title of the item (book, magazine, journal, serial, symposium, catalog, etc.)., date, page(s), volume-issue number(s), publisher, city and/or country where published.	T ²
•	4	Xiang et al., "Fully Quantum Mechanical Energy Optimization for Protein-Ligand Structure," Journal of Computational Chemistry 25(12):1431-1437 (2004)	
	5	Zhang et al., "Full ab initio Computation of Protein-Water Interaction Energies," Journal of Theoretical and Computational Chemistry 3(1):43-49 (2004)	
	6	Zhang et al., "Molecular Caps for Full Quantum Mechanical Computation of Peptide-Water Interaction Energy," Journal of Computational Chemistry 24(15):1846-1852 (2003)	
	7	Zhang et al., "Molecular Fractionation with Conjugate Caps for Full Quantum Mechanical Calculation of Protein-Molecule Interaction Energy," <i>Journal of Chemical Physics</i> 119(7):3599-3605 (2003)	
	8	Zhang et al., "New Advance in Computational Chemistry: Full Quantum Mechanical ab Initio Computation of Streptavidin-Biotin Interaction Energy," J. Phys. Chem. 107:12039-12041 (2003)	
	9	Zhang et al., "Quantum Mechanical Map for Protein-Ligand Binding with Application to β-Trypsin/Benzamidine Complex," <i>Journal of Chemical Physics</i> 120(3):1145-1148 (2004)	
			-

Examiner		Date		
Signature		Considered		

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